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THE Be-Pt (BERYLLIUM-PLATINUM) SYSTEM

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This paper was prepared for submittal to
Bulletin of Alloy Phase Diagrams

November 15, 1985

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Equilibrium Diagram

Available information on the Be-Pt phase diagram is fragmentary as is shown in Fig. 1. Intermediate phases reported in the literature thus far are: (1) Be_{12}Pt (Mn_{12}Th -type); (2) Be_5Pt (AuBe_5 -type); (3) $\text{Be}_{21}\text{Pt}_5$ ($\text{Ni}_5\text{Zn}_{21}$ -type); (4) BePt (CsCl -type); and (5) BePt_{15} (?). Of these, the latter three may need further examination for confirmation. Assessment is based on comparisons with the Be-Pd system [850ka] which shows certain similarities.

Liquidus. No experimental data are available. Assuming that (1) the solid solubility range of (Pt) phase is restricted up to the melting point as in the (Pd) phase in Be-Pd phase diagram [850ka] and (2) the thermodynamic properties of liquid Be-Pt are similar to that of Be-Pd, then the Pt-end of the $L/[L+(\text{Pt})]$ liquidus can be estimated. However, using the similar approach for the $L/[L+(\alpha\text{Be})]$ liquidus is not as reliable because the (αBe) phase may have a sizable solubility range at high temperatures ($>1200^\circ\text{C}$) as can be projected from the solubility data in Table 1.

(βBe) and (αBe) Terminal Solid Solutions. The melting point of (βBe) and the (βBe) \rightarrow (αBe) allotropic transformation temperature are 1289 ± 4 and $1270\pm6^\circ\text{C}$, respectively [85BAP]. [61Pic] reported that Pt may dissolve appreciably in (βBe) but he did not give values. The solubility limit of Pt in (αBe) has been reported as given in Table 1.

(Pt) Terminal Solid Solution. The melting point of Pt is 1769.0°C [Melt]. The reported solubility limit of Be in Pt is listed in Table 2.

Intermediate Phases. The existence of the two Be-rich phases, Be_{12}Pt and Be_5Pt , seem to be well established by XRD studies of single crystals and powders [58Bat] (see Table 3). The latter phase is likely to have some homogeneity range because of

observed variations in lattice parameters (Table 3) [58Bat]. [59Pai] reacted powders of Be_{13}Pt and Be_5Pt at 1400°C in vacuum and obtained ambiguous results, but these may be interpreted as further confirmation for Be_5Pt , based on his finding a cubic phase with $a = 0.6003$ nm. Diffusion couple measurements by [75Mat] suggested that a phase exists with up to 60 wt.% Pt (6.5 at.% Pt) and is stable at 1000 and 1100°C but unstable at 900°C . This phase would appear to correspond to Be_{12}Pt . In addition, [75Mat] found two other distinguishable phases having compositions 85 to 87 wt.% Pt (20.7 to 23.6 at.% Pt) and 93 to 94 wt.% Pt (38 to 42 at.% Pt); the former was believed to correspond to $\text{Be}_{21}\text{Pt}_5$ (19.2 at.% Pt).

The $\text{Be}_{21}\text{Pt}_5$ phase was identified as having a deformed γ brass-type structure by [35Mis] and this was supported by [54Bok], but lattice parameters were not reported in either paper. [75Mat] gave solubility ranges as a function of temperature for this phase as follows:

Temperature, $^\circ\text{C}$	Range, wt.% Pt (at.% Pt)
1100	85 to 87 (20.7 to 23.6)
1000	85 to 86 (20.7 to 22.1)
900	85 (20.7)

Further study here is required.

[59Pai] also reacted an equiatomic powder mixture at 1400°C and found a cubic phase with $a = 0.280$ nm. He compared it to BePd (i.e. B2, CsCl-type) and this agrees with the prediction of structure type and lattice parameter of [80Tan].

Diffusion couple measurements by [75Mat] revealed a phase consisting of 42 at.% Pt at 900°C and 38 at.% Pt at 1000°C . No corresponding phase has been reported in this composition range.

Crystal Structures

Available data on crystal structures are summarized in Table 3.

Thermodynamics

No information is available on the thermodynamic properties of Be-Pt alloys. In order to attempt an estimate of the liquidus near the melting point of Pt, the following assumptions have been made: (1) the (Pt) phase has no solubility range; (2) the lattice stability parameter of solid Pt is given as $-19690 + 9.623T$ J/mol (reference state the liquid), based on the heat of fusion given by [83Cha]; and (3) the Be-Pt liquid behaves almost like Be-Pd liquid [850ka] ($G^{\text{ex}} = -100X(1-X)$ kJ/mol). Fig. 1 shows the calculated liquidus based on these assumptions as well as that from the ideal solution model for comparison. This calculated liquidus is probably not very much different from the actual liquidus near the melting point of Pt.

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* Indicates key paper.

Acknowledgements

Be-Pt evaluation contributed by L.E. Tanner, L-217, Lawrence Livermore National Laboratory, P.O. Box 808, Livermore, CA 94550 and H. Okamoto, B77G, Lawrence Berkeley National Laboratory, Berkeley, CA 94720. Work was supported by the U.S. Department of Energy under contract no. W-7405-Eng-48 and American Society for Metals (ASM). Literature searched through 1984. Part of the bibliographic search was provided by ASM. L.E. Tanner and H. Okamoto are ASM/NBS Data Program Category Editors for binary beryllium alloys.

Table 1 Solubility Limit of Pt in (α Be)

Reference	Composition, at.% Pt	Comments
50Kau 75Mat	<1.1 0.63 at 1200°C	... decreases with decreasing temperature

Table 2 Solubility Limit of Be in (Pt)

Reference	Composition, at.% Be	Comments
41Fro 48Nem 75Mat	1.28 5.15 0.86 to 1.07	... could be 3.15 at.% Be [Hansen] unchanged between 900 and 1100°C

Table 3 Be-Pt Crystal Structure and Lattice Parameter Data

Phase	Composition range, at.% Pt	Pearson symbol	Struktur- bericht designation	Space group	Proto- type	Lattice parameters, nm		Reference
						a	c	
(β Be)....	0 to ?	cI2	A2	Im3m	W	0.25515	...	[King2]
(α Be)....	~0(a)	hP2	A3	P6 ₃ /mmc	Mg	0.22857	0.35839	[King1]
Be ₇ Pt....	7.7	tI26	D2 _b	I4/mmm	Mn ₁₂ Th	0.7237	0.4252	[58Bat]
Be ₅ Pt....	16.7	cF24	C15 _b	F43m or F23	AuBe ₅	0.6004 to 0.5978	...	[58Bat]
Be ₄ Pt ₅ ?..	19.2	Cubic	DB ₁₋₃	?	Ni ₅ Zn ₂₁	?	...	[54Bok]
BePt.....	50	cP2	B2	Pm3m	CsCl	0.280	...	[59Pai, 80Tan]
BePt ₁₅ ?...	93.8	(hexagonal)		?	?	2.213	2.712	[62Bro]
(Pt)....	~100(b)	cF4	A1	Fm3m	Cu	0.39233	...	[King1]

(a) See Table 1.

(b) See Table 2.

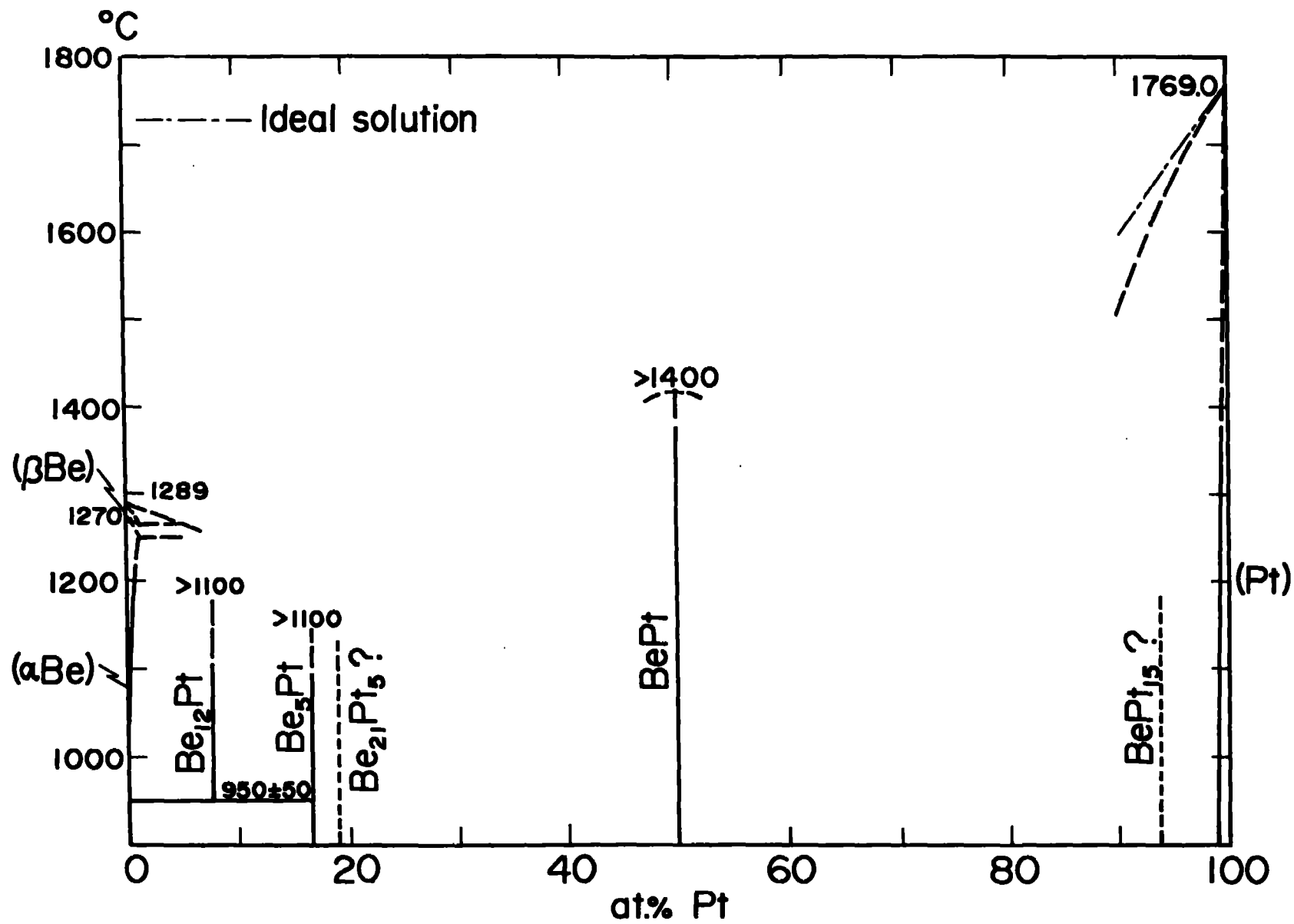


Fig. 1